

Vortex structures of a two-component BEC for large anisotropies

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We calculate the vortex structures of an elongated two-component Bose-Einstein condensate. We study how these structures depend on the intra-component and inter-component interaction strengths. We present analytic and numeric results respectively at weak and strong interactions; finding lattices with different interlocking geometries: triangular, square, rectangular and double-core.

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I. INTRODUCTION

One of the most exciting recent developments in cold atom experiments has been the production of artificial gauge fields, which couple to neutral atoms in the same way that magnetic fields couple to charged particles [1, 2]. While the greatest excitement surrounds the possibility of producing analogs of fractional quantum Hall states or topological insulators [3], these experiments also allow one to perform some extremely interesting experiments on vortices in Bose-Einstein condensates. These structures are particularly rich for multi-component gases [4–7]. A key difference between the latest NIST experiments, and previous ones on rotating condensates [8–13], is that the techniques used to generate the artificial gauge fields naturally lead one to consider very narrow geometries. Here we theoretically investigate the vortex structures in a two-component Bose-Einstein condensate confined to a relatively narrow channel. We find a rich phase diagram which complements our understanding of the isotropic 2D system [4–7].

In the NIST experiments the hyperfine states of ^{87}Rb are coupled by a series of Raman lasers. In the dressed state picture [1, 2], taking the strong Raman field limit, one arrives at an effective single particle Hamiltonian

$$H_1 = [p_x - A_x(y)]^2 / 2m + p_y^2 / 2m + p_z^2 / 2m + V_{\text{eff}}. \quad (1)$$

The effective gauge field is bounded, $-\hbar k_R < A_x(y) < \hbar k_R$, where $\hbar k_R$ is the momentum kick from absorbing a photon from one beam and releasing it into another. One can generate arbitrarily large magnetic fields $B = \partial_y A_x$, but due to the bounded nature of A_x , these can only exist over a finite extent in the y direction. It is therefore natural to choose the effective potential $V_{\text{eff}}(y, z)$ to restrict the particles to a small region of y, z , modeling it as $V_{\text{eff}}(y, z) = m\omega_y^{*2}y^2/2 + m\omega_z^2z^2/2$. V_{eff} includes both the external potential and one induced by the Raman lasers, and ω_y^* is an effective trap frequency.

The impact of the magnetic field on a single component Bose-Einstein condensate in such a narrow channel has been extensively studied [17–19]. As one increases the magnetic field from zero, the Bogoliubov spectrum

starts to develop a “roton” minimum, whose energy falls with increasing B . When the roton energy hits zero, the system becomes unstable to undulations. At higher magnetic fields vortices enter the channel. Here we extend these results to the two-component gas.

The vortex structures in isotropic 2D two-component condensates are quite rich [4, 5]. Depending on the relative strength of the inter-component and intra-component scattering one can find interpenetrating lattices with different geometries: triangular, square, rectangular and double-core. Additionally, one sees interesting hysteresis [5]. We find similar results in the anisotropic geometry.

A final motivation for thinking about the role of magnetic fields in very narrow geometries comes from solid state physics. In 2005, Seidel et al [16] showed that the quantum Hall effect in narrow tori can be connected to charge density waves, with the fractionally charged vortices mapping onto fractionally charged domain walls. In this paper we will focus on the Bose condensed regime, and will not be able to comment on this interesting physics. We do note, however, that our system undergoes a charge density wave instability before vortices enter the system. Investigating this instability in the low density “quantum” limit may be fruitful, even though the physics is likely to be very different from that in [16].

II. MODEL

There are four important lengths in Eq. (1), $d_z = \sqrt{\hbar/m\omega_z}$, $d_y = \sqrt{\hbar/m\omega_y^*}$, $\ell = \sqrt{\hbar/B}$, and $R = \hbar k_R/\bar{B}$, where \bar{B} is the peak value of B . These correspond to the trap length in the z and y direction, the magnetic length, and the spatial range over which the effective magnetic is nonzero. If d_z is sufficiently small, the extent in z direction is sufficiently compressed and the 3D gases can be treated as 2D cloud. As we argue below, a sufficient condition will be $d_z \lesssim \ell$. If $d_y \ll R$, the cloud will be confined to a region where the vector potential varies linearly and we can approximate: $A_x = \bar{B}y$. In terms of the raw experimental parameters, $\bar{B} = \delta'\hbar k_R/\Omega$, where

Ω is the Raman Rabi frequency, and δ' is the gradient of two-photon detuning along y direction.

Taking $\Psi_1(\mathbf{r})$ and $\Psi_2(\mathbf{r})$ to annihilate atoms in the two (pseudo)-spin states, the short range interactions will be

$$H_{\text{int}} = \frac{1}{2} \int d\mathbf{r} (g_1 \Psi_1^\dagger \Psi_1^\dagger \Psi_1 \Psi_1 + g_2 \Psi_2^\dagger \Psi_2^\dagger \Psi_2 \Psi_2 + 2g_{12} \Psi_1^\dagger \Psi_2^\dagger \Psi_2 \Psi_1). \quad (2)$$

The coupling constants are related to scattering lengths a_s via $g = 4\pi\hbar^2 a_s/m$. We will consider the case $g_1 = g_2 = g$, and $g_{12} = g_m$. For most experiments $g_1 \approx g_2 \approx g_{12}$, but one gains insight by relaxing this condition.

Following Sinha and Shlyapnikov [17], we diagonalize the single particle Hamiltonian in Eq. (1). The eigenstates are labeled by three quantum numbers K, n, n' , with energies

$$E_{nn'}(K) = \mathcal{E}K^2 + n\hbar\omega_z + n'\hbar\tilde{\omega}_c. \quad (3)$$

where $\tilde{\omega}_c^2 = \omega_y^{*2} + \omega_c^2$, $\omega_c = \bar{B}/m$, and $\mathcal{E} = \hbar^2\omega_y^{*2}/4m\tilde{\omega}_c^2\tilde{\ell}^2$, $\tilde{\ell} = \sqrt{\hbar/m\tilde{\omega}_c}$, and we have neglected the zero-point energy. The continuous variable $K = \sqrt{2}\tilde{\ell}k$ is a dimensionless label of the momentum k along the x direction, while n and n' are discrete quantum numbers corresponding to the number of nodes in the z and y directions. If the interaction energy per particle $\langle H_{\text{int}}/N \rangle$ and the characteristic “kinetic energy” $\langle \mathcal{E}K^2 \rangle$ are small compared to $\hbar\tilde{\omega}_c$ and $\hbar\omega_z$ one can truncate to the single particle eigenstates with $n = n' = 0$, which are of the form

$$\phi_K(x, y, z) = \frac{1}{\sqrt{\pi\tilde{\ell}d_z L}} e^{i\frac{Kx}{\sqrt{2}\tilde{\ell}}} e^{-\frac{(y-y_K)^2}{2\tilde{\ell}^2}} e^{-\frac{z^2}{2d_z^2}} \quad (4)$$

where $y_K = \sqrt{2}\omega_c K \tilde{\ell} / 2\tilde{\omega}_c$ and L is the length of x direction.

By taking the system sufficiently dilute, it is easy to arrange $\langle H_{\text{int}}/N \rangle \ll \hbar\tilde{\omega}_c \sim \hbar\omega_z$. The other condition, $\mathcal{E}K^2 \ll \hbar\tilde{\omega}_c$, will be valid at strong magnetic fields. For our vortex lattices, we find the characteristic dimensionless wave-number to be $K \sim 1$, thus requiring $\omega_y^* \ll \omega_c$. Combining this with the previous constraint, $\omega_y^* \gg m\omega_c^2/\hbar k_R^2$, we see that our approximations break down unless the magnetic length is much larger than the wavelength of the Raman lasers, ie. $\omega_c \ll \hbar k_R^2/m$. This also establishes our requirement $d_z \lesssim \ell$.

Letting a_K annihilate the state in Eq. (4), the N -body Hamiltonian is

$$H/\mathcal{E} = \sum_K K^2 [a_{K\uparrow}^\dagger a_{K\uparrow} + a_{K\downarrow}^\dagger a_{K\downarrow}] + \frac{\beta}{N} \sum_q F_\uparrow^\dagger(q) F_\uparrow(q) + \frac{\beta}{N} \sum_q F_\downarrow^\dagger(q) F_\downarrow(q) + \frac{2\beta_m}{N} \sum_q F_{\uparrow\downarrow}^\dagger(q) F_{\uparrow\downarrow}(q) \quad (5)$$

where

$$F_{\sigma\tau}(q) = \sum_{K_1 K_2} \delta_{q-K_1-K_2} e^{-\frac{1}{8}(K_1-K_2)^2} a_{K_1\sigma} a_{K_2\tau}. \quad (6)$$

We take the continuum limit $\sum_K \rightarrow (\sqrt{2}L/4\pi\tilde{\ell}) \int dK$, and $\delta_K \rightarrow (2\sqrt{2}\pi\tilde{\ell}/L)\delta(K)$. The effective 1D parameters are $\beta = Nmg\tilde{\omega}_c^2\tilde{\ell}/\pi L d_z \omega_y^{*2} \hbar^2$, $\beta_m = \beta g_m/g$. From these definitions of the effective interaction parameters β, β_m , we see that increasing g, g_m has the same effect as increasing the magnetic field B , increasing the confinement ω_z , or reducing the confinement ω_y^* .

We consider a variational wavefunction corresponding to a Bose-Einstein condensate which is periodic in the x direction, with dimensionless wavelength $2\pi/K_0$, corresponding to a physical wavelength $\lambda = 2\sqrt{2}\pi\tilde{\ell}/K_0$

$$|\psi\rangle = \exp \left[-\frac{N}{2} + \sum_{\sigma} \sqrt{N_{\sigma}} \left(\sum_n C_{n\sigma} a_{nK_0\sigma}^\dagger \right) \right] |\text{vac}\rangle. \quad (7)$$

Where $\sigma = \uparrow, \downarrow$ for two-component BEC. N_{σ} is the number of particles in state σ , and $N = N_{\uparrow} + N_{\downarrow}$. The coefficients $C_{n\sigma}$ are normalized to $\sum_n |C_{n\sigma}|^2 = 1$. In place of such a coherent state ansatz, some authors prefer to work with a “Fock” state

$$|\psi\rangle_F = \frac{1}{\sqrt{N_{\uparrow}! N_{\downarrow}!}} \left(\sum_n C_{n\uparrow} a_{nK_0\uparrow}^\dagger \right)^{N_{\uparrow}} \left(\sum_n C_{n\downarrow} a_{nK_0\downarrow}^\dagger \right)^{N_{\downarrow}} |\text{vac}\rangle. \quad (8)$$

For the quantities we are interested in, $|\psi\rangle$ and $|\psi\rangle_F$ are equivalent, and the variational parameters $C_{n\sigma}$ have the same meaning in each case: $|C_{n\sigma}|^2$ is the fraction of spin σ particles with momentum nK_0 .

We take $N_{\uparrow} = N_{\downarrow}$, and further restrict ourselves to considering symmetric or antisymmetric wavefunctions: $C_{n,\sigma} = \pm C_{-n,\sigma}$. We classify our state by the value of K_0 and ξ the number of non-zero C_n ’s which are needed to minimize the energy. For example, as illustrated in Fig. 1, for $(\beta, \beta_m) \sim (3, 1)$, we need only one n for each component: $n = 0$. We refer to this state as $(\xi_{\uparrow}, \xi_{\downarrow}) = (1, 1)$. This should be contrasted with the case at $(\beta, \beta_m) \sim (5, 2.5)$, where the energy is minimized by taking $n = -1, 0, 1$ in both components, a state we label as $(\xi_{\uparrow}, \xi_{\downarrow}) = (3, 3)$. Analytic expressions for the energies with small numbers of components are given in appendix A.

III. RESULTS AT SMALL β, β_m

The number of expansion parameters, $(\xi_{\uparrow}, \xi_{\downarrow})$, grow with the magnitude of β, β_m . For small β and β_m , we only need a small number of terms in our wavefunctions, and we may use the analytic expressions in Eq. A1-A15 to find the lowest energy state. We obtain a series of phase, as demonstrated in Fig. 1(a).

At $\beta_m = 0$, the \uparrow and \downarrow atoms decouple, and the physics is identical to the single component case, $\xi_{\uparrow} = \xi_{\downarrow} = \xi$. When $\beta < 9.8$, the ground state wavefunction has only one term ($\xi = 1$), which displays a Gaussian

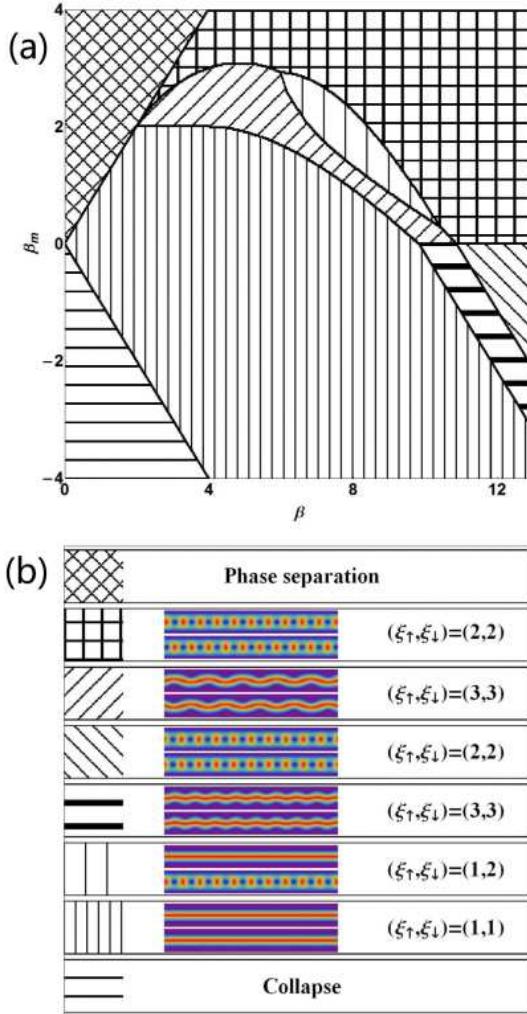


FIG. 1: Top panel: phase diagram at small dimensionless intra-species and inter-species interactions β, β_m . Larger β, β_m corresponds to larger g, g_m , larger magnetic field, larger confinement ω_z , or weaker confinement ω_y^* . Hatched patterns represent states described by different number of Fourier components in each spin state: $(\xi_\uparrow, \xi_\downarrow)$. Bottom panel: the density profiles of the two-component wavefunction in the corresponding regimes. A color key for the density patterns is shown in Fig. 2.

shape along y direction. As β increases past $\beta = 9.8$, the $\xi = 1$ wavefunction becomes unstable, and it undergoes a second-order phase transition to a density wave with $\xi = 3$. As β increases to $\beta = 10.8$, a first-order phase transition occurs to a state with $\xi = 2$, characterized by a single row of vortices. These results have been extensively studied in Ref. [19].

In the regime of repulsive inter-component interaction ($\beta_m > 0$), the \uparrow and \downarrow particles try to avoid each-other. For strong repulsive interaction ($\beta_m > \beta$), the two components undergo a microscale phase separate [14], which needs large (presumably infinite) xi to describe. For weak repulsive interaction ($\beta_m < \beta$) rich structures, illustrated in Fig. 1 develop. For some β, β_m , we find

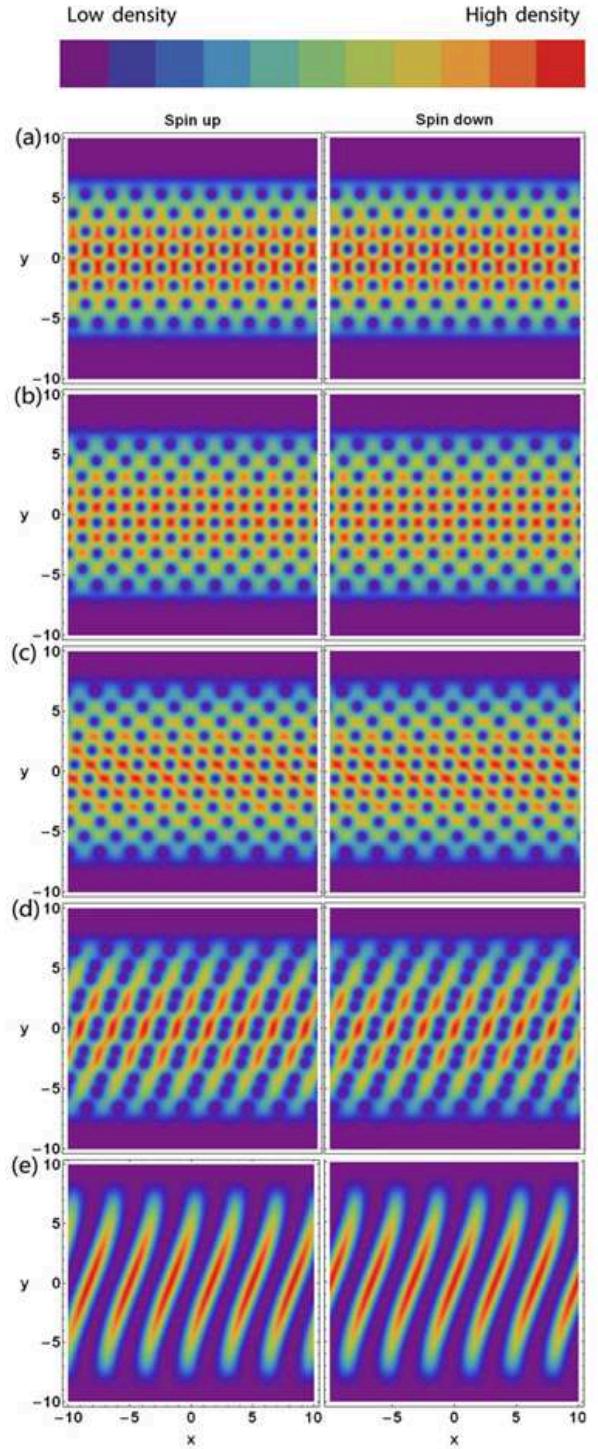


FIG. 2: The density profiles of two-component BEC at large β, β_m , where x, y are in the unit of $\sqrt{2\ell}$. For (a) – (e), the parameters are $\beta_m/\beta = 0.1, 0.5, 0.8, 1, 1.5$ respectively, with $\beta = 1000$. A color key is shown at the top.

$\xi_\uparrow \neq \xi_\downarrow$. For example, at $(\beta, \beta_m) = (8, 2)$, $\xi_\uparrow = 1$ and $\xi_\downarrow = 2$, and the \downarrow atoms have a row of vortices, while the \uparrow atoms show no structure. Under these circumstances

there is a degenerate state with the \uparrow and \downarrow wavefunctions reversed. For other β, β_m , there is a symmetry between the two components with $\xi_\uparrow = \xi_\downarrow$. The vortices or density corrugations are displaced by half a period so that the density maxima of the \uparrow atoms line up with the density minima of the others. For example, when $(\beta, \beta_m) = (10, 3)$, each component displays a single row vortices, and the wavefunctions are related by a translation.

In the regime of attractive inter-component interaction ($\beta_m < 0$), one wants to maximize the overlap of the \uparrow and \downarrow wavefunctions. Generically, this means that the wavefunction of each component is identical, and the problem reduces to the single component case, but with a renormalized interaction $\beta \rightarrow \beta + \beta_m$. The phase diagram for $\beta_m < 0$ can be calculated from the phase diagram at $\beta_m = 0$ by mapping each point: $(\beta, \beta_m) \rightarrow (\beta - \beta_m, 0)$. For strong attractive interaction ($-\beta_m > \beta$), the BEC is unstable and expected to collapse, similar to the case of a single component BEC with attractive interaction [15].

IV. RESULTS AT LARGE β, β_m

For large β and β_m , the analytic expressions for the energy become unwieldy. We numerically minimize the expectation $\langle H \rangle$, varying $\{C_{n\sigma}, K_0\}$ in Eq. (7). The energy landscape has many local minima, and we use a range of starting parameters to try to find the absolute minimum. We cannot rule out the existence of even lower energy states. Moreover, for some parameters we found that the energy differences between competing minima became extremely small. In an experiment it is doubtful that one would find the true minimum energy state. Rather than systematically exploring the large β physics, we simply show a few examples.

Our results, illustrated in Fig. 2, are richer than those seen in the single component gas [19], showing structures similar to those in isotropic 2D studies [4, 5]. When $\beta_m \ll \beta$ one finds two interlocking triangular lattices, as in Fig. 2(a) where $\beta_m/\beta = 0.1$. As one

increases β_m the lattice structure changes: interlocking square/rectangular lattices are shown in Fig. 2(b,c) where $\beta_m/\beta = 0.5, 0.8$. When $\beta_m > \beta$ the non-rotating system would be expected to phase separate. Here, one finds more intricate vortex structures at $\beta_m \sim \beta$. At $\beta_m/\beta = 1$, we find double-core vortices (cf. [5]), as in Fig. 2(d). At $\beta_m/\beta = 1.5$ we find stripes, which are a microscopic version of phase separation. Close inspection of the image in Fig. 2(e), shows vortex cores in the low density regions.

V. SUMMARY AND CONCLUSIONS

We have investigated the vortex structures in a two-component BEC with an artificial magnetic field, in an elongated geometry. Compared to the single component gas, the two-component vortex structures are more intricate.

To experimentally investigate these structures, one needs to find a system where the interspecies interactions can be tuned relative to the intraspecies. One promising approach is to use different atomic species for the two (pseudo)-spin states, and take advantage of an interspecies Feshbach resonance [20]. Some of the other newly condensed atomic systems may also be favorable [21]. If one cannot separately tune β and β_m , one can still change the magnitude of them, fixing β_m/β . As seen in Fig. 1, such a cut through the phase diagram can still be quite rich, especially if $\beta_m/\beta \sim 0.25$.

VI. ACKNOWLEDGMENT

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Appendix A: Analytic results for small number of components.

Here we give analytic results for the energies of the states defined in Eq. (7), truncating the n -sums for each σ , and assuming symmetry/antisymmetry about the origin for each component.

State $(\xi_\uparrow, \xi_\downarrow) = (1, 1)$ is unique, and has energy

$$E_{1,1} = \frac{1}{2}(\beta + \beta_m) \quad (\text{A1})$$

State $(\xi_\uparrow, \xi_\downarrow) = (1, 2)$, with

$$\psi_\uparrow = \phi_{0\uparrow} \quad (\text{A2})$$

$$\psi_\downarrow = \frac{\sqrt{2}}{2}(\phi_{K_0\downarrow} + \phi_{-K_0\downarrow}) \quad (\text{A3})$$

is unique up to translation, and has energy

$$E_{1,2} = \frac{1}{2}K_0^2 + \frac{1}{4}\beta(e^{-K_0^2} + \frac{3}{2}) + \frac{1}{2}\beta_m e^{-\frac{1}{4}K_0^2} \quad (\text{A4})$$

State $(\xi_\uparrow, \xi_\downarrow) = (1, 3)$, with

$$\psi_\uparrow = \phi_{0\uparrow} \quad (\text{A5})$$

$$\psi_\downarrow = \sqrt{1 - 2|\epsilon|^2}\phi_{0\downarrow} - \epsilon(\phi_{K_0\downarrow} + \phi_{-K_0\downarrow}) \quad (\text{A6})$$

has energy

$$\begin{aligned} E_{1,3} &= \frac{1}{2}(\beta + \beta_m) \\ &+ [K_0^2 + \beta(2e^{-\frac{1}{4}K_0^2} - e^{-\frac{1}{2}K_0^2} - 1) + \beta_m(e^{-\frac{1}{4}K_0^2} - 1)]\epsilon^2 \\ &+ \beta \left(e^{-K_0^2} + 2e^{-\frac{1}{2}K_0^2} - 4e^{-\frac{1}{4}K_0^2} + \frac{3}{2} \right) \epsilon^4 \end{aligned} \quad (\text{A7})$$

For state $(\xi_\uparrow, \xi_\downarrow) = (2, 2)$, there are two distinct extremal states. Both components could be symmetric about the origin ($C_{1\uparrow} = C_{-1\uparrow} = C_{1\downarrow} = C_{-1\downarrow} = 1/\sqrt{2}$), or one of them could be antisymmetric ($C_{1\uparrow} = C_{-1\uparrow} = C_{1\downarrow} = -C_{-1\downarrow} = 1/\sqrt{2}$). This gives energies

$$E_{2,2}^s = K_0^2 + \frac{1}{4}(\beta + \beta_m) + \frac{1}{2}(\beta + \beta_m)e^{-K_0^2} \quad (\text{A8})$$

$$E_{2,2}^a = E_{2,2}^s - \frac{1}{2}\beta_m e^{-K_0^2} \quad (\text{A9})$$

For state $(\xi_\uparrow, \xi_\downarrow) = (2, 3)$, the \uparrow component can be symmetric or antisymmetric, resulting in energies

$$\begin{aligned} E_{2,3}^s &= \frac{1}{2}K_0^2 + \frac{1}{4}\beta(e^{-K_0^2} + \frac{3}{2}) + \frac{1}{2}\beta_m e^{-\frac{1}{4}K_0^2} \\ &+ [K_0^2 + \beta(2e^{-\frac{1}{4}K_0^2} - e^{-\frac{1}{2}K_0^2} - 1)]\epsilon^2 \\ &+ \frac{1}{2}\beta_m(1 - 2e^{-\frac{1}{4}K_0^2} + 2e^{-K_0^2})\epsilon^2 \\ &+ \beta(\frac{3}{2} - 4e^{-\frac{1}{4}K_0^2} + 2e^{-\frac{1}{2}K_0^2} + e^{-K_0^2})\epsilon^4 \end{aligned} \quad (\text{A10})$$

$$E_{2,3}^a = E_{2,3}^s - \beta_m e^{-K_0^2}\epsilon^2 \quad (\text{A11})$$

For state $(\xi_\uparrow, \xi_\downarrow) = (3, 3)$, the optimal wavefunction is

$$\psi_\uparrow = \sqrt{1 - 2|\epsilon|^2}\phi_{0\uparrow} + i\epsilon(\phi_{K_0\uparrow} + \phi_{-K_0\uparrow}) \quad (\text{A12})$$

$$\psi_\downarrow = \sqrt{1 - 2|\epsilon|^2}\phi_{0\downarrow} \pm i\epsilon(\phi_{K_0\downarrow} + \phi_{-K_0\downarrow}) \quad (\text{A13})$$

with ϵ real.

If $\beta > \beta_m > 0$, the negative sign has lower energy,

$$\begin{aligned} E_{3,3}^- &= \frac{1}{2}(\beta + \beta_m) \\ &+ [2K_0^2 + \beta(-2e^{-\frac{1}{2}K_0^2} + 4e^{-\frac{1}{4}K_0^2} - 2)]\epsilon^2 \\ &+ \beta_m(2e^{-\frac{1}{2}K_0^2} - 2)\epsilon^2 \\ &+ \beta(2e^{-K_0^2} + 4e^{-\frac{1}{2}K_0^2} - 8e^{-\frac{1}{4}K_0^2} + 3)\epsilon^4 \\ &+ \beta_m(2e^{-K_0^2} - 4e^{-\frac{1}{2}K_0^2} + 3)\epsilon^4 \end{aligned} \quad (\text{A14})$$

If $\beta > 0, \beta_m < 0$, the positive sign has lower energy,

$$E_{3,3}^+ = E_{3,3}^- + 4\beta_m(e^{-K_0^2/4} - e^{-K_0^2/2})(\epsilon^2 + 2\epsilon^4) \quad (\text{A15})$$